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NEWS 27 JUN 13 USPATFULL and USPAT2 updated with 11-character patent numbers for U.S. applications
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NEWS 32 JUN 30 STN on the Web enhanced with new STN AnaVist
Assistant and BLAST plug-in

NEWS 33 JUN 30 STN AnaVist enhanced with database content from EPFULL

NEWS EXPRESS JUNE 27 08 CURRENT WINDOWS VERSION IS V8.3,
AND CURRENT DISCOVER FILE IS DATED 23 JUNE 2008.

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NEWS LOGIN Welcome Banner and News Items

NEWS IPC8 For general information regarding STN implementation of IPC 8

Enter NEWS followed by the item number or name to see news on that specific topic.

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FILE 'HOME' ENTERED AT 10:43:19 ON 24 JUL 2008

=> file reg
COST IN U.S. DOLLARS
SINCE FILE ENTRY SESSION
0.21 0.21
FULL ESTIMATED COST

FILE 'REGISTRY' ENTERED AT 10:43:24 ON 24 JUL 2008
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STRUCTURE FILE UPDATES: 22 JUL 2008 HIGHEST RN 1035393-16-8
DICTIONARY FILE UPDATES: 22 JUL 2008 HIGHEST RN 1035393-16-8

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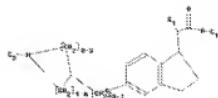
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=>
Uploading C:\Program Files\STNEXP\Queries\10537630_4.str



chain nodes :
 10 11 12 13 16 17 20 21 29
 ring nodes :
 1 2 3 4 5 6 7 8 9 25 26 27 28
 chain bonds :
 2-20 7-10 10-11 10-17 11-12 11-13 13-16 20-21 21-25 28-29
 ring bonds :
 1-2 1-6 2-3 3-4 4-5 5-6 5-7 6-9 7-8 8-9 25-26 25-27 26-28 27-28
 exact/norm bonds :
 2-20 10-17 11-12 11-13 13-16 20-21 28-29
 exact bonds :
 5-7 6-9 7-8 7-10 8-9 10-11 21-25 25-26 25-27 26-28 27-28
 normalized bonds :
 1-2 1-6 2-3 3-4 4-5 5-6
 isolated ring systems :
 containing 1 : 25 :

G1:H,Cb,Ak

G2:O,S,SO2

G3:Cb,Cy,Hy

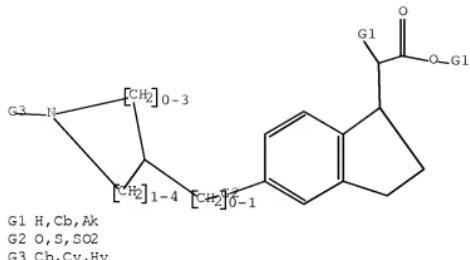
Match level :
 1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:CLASS
 11:CLASS 12:CLASS 13:CLASS 16:CLASS 17:CLASS 20:CLASS 21:CLASS 25:Atom
 26:Atom 27:Atom 28:Atom
 29:CLASS

L1 STRUCTURE UPLOADED

=> d L1

L1 HAS NO ANSWERS

L1 STR



Structure attributes must be viewed using STN Express query preparation.

=> file caplus		
COST IN U.S. DOLLARS	SINCE FILE	TOTAL
FULL ESTIMATED COST	ENTRY	SESSION
	0.46	0.67

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FILE COVERS 1907 - 24 Jul 2008 VOL 149 ISS 4
 FILE LAST UPDATED: 23 Jul 2008 (20080723/ED)

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=> s L1 SSS full
 REGISTRY INITIATED
 Substance data SEARCH and crossover from CAS REGISTRY in progress...
 Use DISPLAY HITSTR (or FHITSTR) to directly view retrieved structures.

FULL SEARCH INITIATED 10:43:52 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 7928 TO ITERATE

100.0% PROCESSED 7928 ITERATIONS
SEARCH TIME: 00.00.01

8 ANSWERS

L2 8 SEA SSS FUL L1

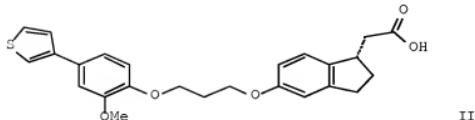
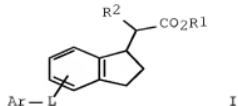
L3 1 L2

=> d ibib abs hitstr 1-
YOU HAVE REQUESTED DATA FROM 1 ANSWERS - CONTINUE? Y/(N):y

L3 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2008 ACS on STN
ACCESSION NUMBER: 2004:565052 CAPLUS Full-text
DOCUMENT NUMBER: 141:123483
TITLE: Preparation of indaneacetic acid derivatives and their
use as pharmaceutical agents
INVENTOR(S): Cantin, Louis-David; Choi, Soongyu; Clark, Roger B.;
Hentemann, Martin F.; Ma, Xin; Rudolph, Joachim;
Liang, Sidney X.; Akuche, Christiana; Lavoie, Rico C.;
Chen, Libing; Majumdar, Duyti; Wickens, Philip L.
PATENT ASSIGNEE(S): Bayer Pharmaceuticals Corporation, USA
SOURCE: PCT Int. Appl., 230 pp.
CODEN: PIXXD2
DOCUMENT TYPE: Patent
LANGUAGE: English
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2004058174	A2	20040715	WO 2003-US40842	20031219
WO 2004058174	A3	20041202		
W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BW, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DC, EC, EE, EG, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NI, NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SY, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW RW: BW, GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG				
CA 2510793	A1	20040715	CA 2003-2510793	20031219
AU 2003299790	A1	20040722	AU 2003-299790	20031219
EP 1578715	A2	20050928	EP 2003-800063	20031219
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, MK, CY, AL, TR, BG, CZ, EE, HU, SK				
JP 2006516251	T	20060629	JP 2004-563903	20031219
US 20060084680	A1	20060420	US 2005-537630	20050603
PRIORITY APPLN. INFO.:			US 2002-435310P	P 20021220
			WO 2003-US40842	W 20031219

OTHER SOURCE(S): MARPAT 141:123483



AB The title compds. [I; R1, R2 = H, alkyl, cycloalkyl; L = (CH2)mX, Y(CH2)nY, etc.; X = O, S, SO, SO2, Y = O, S, SO, SO2, (un)substituted NH; m = 1-3; n = 2-4; Ar = (un)substituted Ph, 5-6 membered heteroaryl containing up to there N atoms] which are useful in the treatment of diseases such as diabetes, obesity, hyperlipidemia, and atherosclerotic diseases, were prepared and formulated. Thus, coupling Et [(1S)-5-[3-(4-bromo-2-methoxyphenoxy)propoxy]-2,3-dihydro-1H-inden-1-yl]acetate (preparation given) with 3-thiopheneboronic acid in the presence of PdCl2(dppf).CH2Cl2, NaHCO3 in DME/H2O followed by treatment of the resulting ester with LiOH afforded (1S)-II.

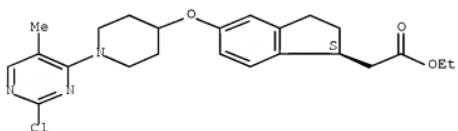
IT 724470-54-6P

RL: PAC (Pharmacological activity); RCT (Reactant); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); RACT (Reactant or reagent); USES (Uses)
(preparation of indaneacetic acid derivs. for treating diabetes, obesity, hyperlipidemia, and atherosclerotic diseases)

RN 724470-54-6 CAPLUS

CN 1H-Indene-1-acetic acid, 5-[(1-(2-chloro-5-methyl-4-pyrimidinyl)-4-piperidinyl)oxy]-2,3-dihydro-, ethyl ester, (1S)- (CA INDEX NAME)

Absolute stereochemistry.



IT 724470-55-7P 724470-56-8P 724470-57-9P
724470-58-0P 724470-59-1P 724470-60-4P
724470-61-5P

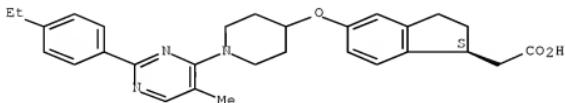
RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(preparation of indaneacetic acid derivs. for treating diabetes, obesity, hyperlipidemia, and atherosclerotic diseases)

RN 724470-55-7 CAPLUS

CN 1H-Indene-1-acetic acid, 5-[(1-[2-(4-ethylphenyl)-5-methyl-4-pyrimidinyl]-4-piperidinyl)oxy]-2,3-dihydro-, (1S)- (CA INDEX NAME)

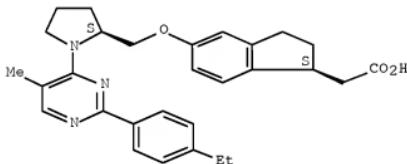
Absolute stereochemistry.



RN 724470-56-8 CAPLUS

CN 1H-Indene-1-acetic acid, 5-[(2S)-1-[2-(4-ethylphenyl)-5-methyl-4-pyrimidinyl]-2-pyrrolidinyl]methoxy]-2,3-dihydro-, (1S)- (CA INDEX NAME)

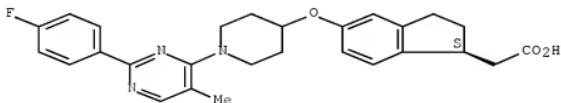
Absolute stereochemistry.



RN 724470-57-9 CAPLUS

CN 1H-Indene-1-acetic acid, 5-[(1-[2-(4-fluorophenyl)-5-methyl-4-pyrimidinyl]-4-piperidinyl)oxy]-2,3-dihydro-, (1S)- (CA INDEX NAME)

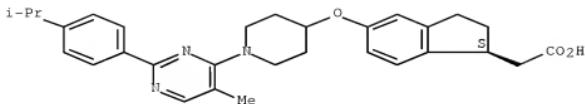
Absolute stereochemistry.



RN 724470-58-0 CAPLUS

CN 1H-Indene-1-acetic acid, 2,3-dihydro-5-[(1-[5-methyl-2-[4-(1-methylethyl)phenyl]-4-pyrimidinyl]-4-piperidinyl)oxy]-, (1S)- (CA INDEX NAME)

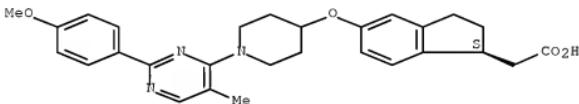
Absolute stereochemistry.



RN 724470-59-1 CAPLUS

CN 1H-Indene-1-acetic acid, 2,3-dihydro-5-[(1-[2-(4-methoxyphenyl)-5-methyl-4-pyrimidinyl]4-piperidinyl)oxy]-, (1S)- (CA INDEX NAME)

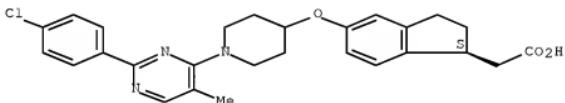
Absolute stereochemistry.



RN 724470-60-4 CAPLUS

CN 1H-Indene-1-acetic acid, 5-[(1-[2-(4-chlorophenyl)-5-methyl-4-pyrimidinyl]4-piperidinyl)oxy]-2,3-dihydro-, (1S)- (CA INDEX NAME)

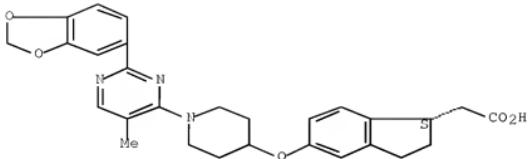
Absolute stereochemistry.



RN 724470-61-5 CAPLUS

CN 1H-Indene-1-acetic acid, 5-[(1-[2-(1,3-benzodioxol-5-yl)-5-methyl-4-pyrimidinyl]4-piperidinyl)oxy]-2,3-dihydro-, (1S)- (CA INDEX NAME)

Absolute stereochemistry.



=> log off

ALL L# QUERIES AND ANSWER SETS ARE DELETED AT LOGOFF

LOGOFF? (Y)/N/HOLD:y

STN INTERNATIONAL LOGOFF AT 10:44:33 ON 24 JUL 2008